Full wwPDB X-ray Structure Validation Report (i)

Aug 16, 2023 – 10:42 AM EDT

PDB ID : 2AOX

Title : Histamine Methyltransferase (Primary Variant T105) Complexed with the Acetylcholinesterase Inhibitor and Altzheimer's Disease Drug Tacrine

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Deposited on : 2005-08-14

Resolution : 3.12 Å(reported)

The following versions of software and data (see references (i)) were used in the production of this report:

|		
MolProbity	:	4.02b-467
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	NOT EXECUTED
EDS	:	NOT EXECUTED
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.35

1 Overall quality at a glance ⓘ

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

The reported resolution of this entry is 3.12 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.
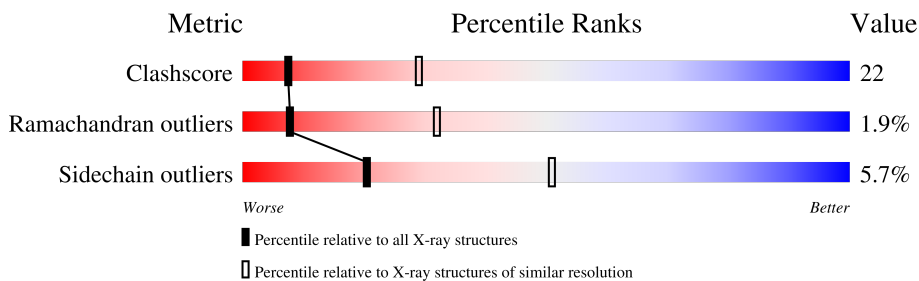
Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
Clashscore	141614	1389 (3.14-3.10)
Ramachandran outliers	138981	1337 (3.14-3.10)
Sidechain outliers	138945	1337 (3.14-3.10)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments of the lower bar indicate the fraction of residues that contain outliers for >=3, 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions <=5%

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	A	292	58% 37%
1	B	292	58% 37%

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

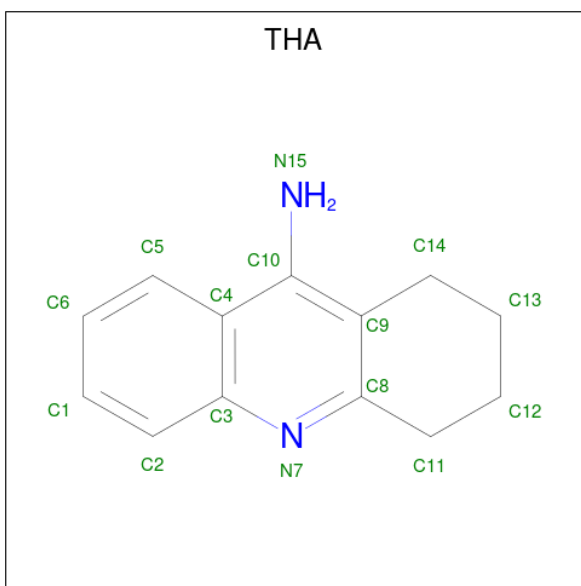| Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
2	THA	A	400	-	X	-	-
2	THA	B	401	-	X	-	-

2   Entry composition ⓘ

There are 3 unique types of molecules in this entry. The entry contains 4511 atoms, of which 0 are hydrogens and 0 are deuteriums.

In the tables below, the ZeroOcc column contains the number of atoms modelled with zero occupancy, the AltConf column contains the number of residues with at least one atom in alternate conformation and the Trace column contains the number of residues modelled with at most 2 atoms.

- Molecule 1 is a protein called Histamine N-methyltransferase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	288	Total	C	N	O	S	0	0	0
			2236	1431	360	430	15			
1	B	288	Total	C	N	O	S	0	0	0
			2236	1431	360	430	15			

- Molecule 2 is TACRINE (three-letter code: THA) (formula: $C_{13}H_{14}N_2$).

THA

Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	C	N	0	0
			15	13	2		
2	B	1	Total	C	N	0	0
			15	13	2		

- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	7	Total 7	O 7	0	0
3	B	2	Total 2	O 2	0	0

3 Residue-property plots ⓘ

These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

Note EDS was not executed.

• Molecule 1: Histamine N-methyltransferase

Chain A:
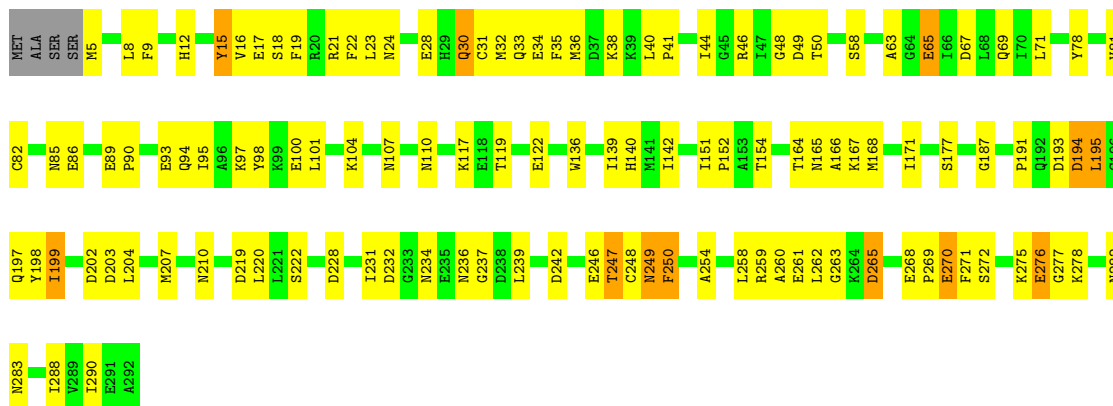
• Molecule 1: Histamine N-methyltransferase

Chain B:
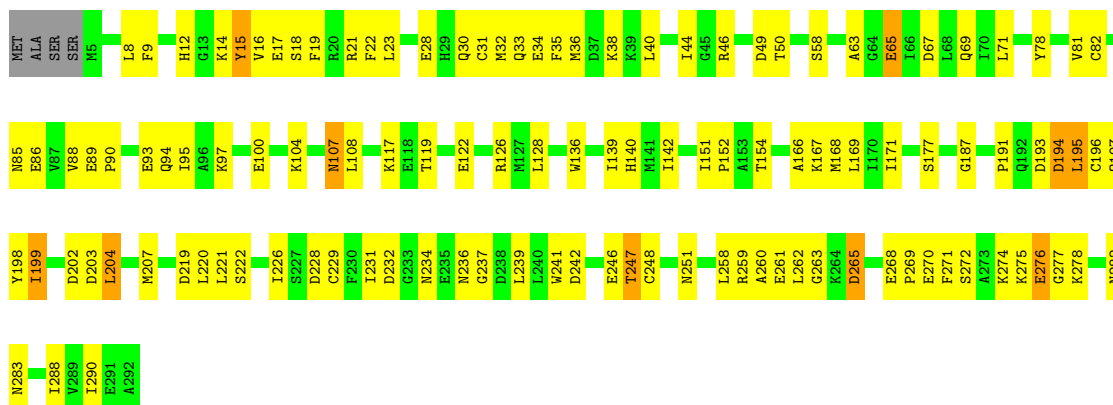
4 Data and refinement statistics ⓘ

Xtriage (Phenix) and EDS were not executed - this section is therefore incomplete.

Property	Value	Source
Space group	P 6	Depositor
Cell constants a, b, c, α, β, γ	132.34Å 132.34Å 64.58Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	25.00 – 3.12	Depositor
% Data completeness (in resolution range)	(Not available) (25.00-3.12)	Depositor
R_{merge}	0.11	Depositor
R_{sym}	(Not available)	Depositor
Refinement program	CNS	Depositor
R, R_{free}	0.229 , 0.285	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	4511	wwPDB-VP
Average B, all atoms (Å^2)	45.0	wwPDB-VP

5   Model quality ⓘ

5.1   Standard geometry ⓘ

Bond lengths and bond angles in the following residue types are not validated in this section: THA

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 5$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	#$\|Z\|$ >5	RMSZ	#$\|Z\|$ >5
1	A	0.50	0/2284	0.65	0/3092
1	B	0.51	0/2284	0.64	0/3092
All	All	0.50	0/4568	0.65	0/6184

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

5.2   Too-close contacts ⓘ

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry-related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	2236	0	2131	95	1
1	B	2236	0	2131	100	0
2	A	15	0	14	0	0
2	B	15	0	14	0	0
3	A	7	0	0	0	0
3	B	2	0	0	0	0
All	All	4511	0	4290	194	1

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 22.

All (194) close contacts within the same asymmetric unit are listed below, sorted by their clash

magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:268:GLU:HG3	1:B:269:PRO:HD2	1.42	1.00
1:A:234:ASN:HD22	1:A:237:GLY:H	1.02	0.94
1:A:268:GLU:HG3	1:A:269:PRO:HD2	1.49	0.94
1:B:142:ILE:HG23	1:B:171:ILE:HD12	1.49	0.93
1:B:234:ASN:HD22	1:B:237:GLY:H	1.00	0.93
1:A:198:TYR:O	1:A:199:ILE:HD12	1.70	0.91
1:B:191:PRO:HG3	1:B:247:THR:HG23	1.55	0.89
1:B:198:TYR:O	1:B:199:ILE:HD12	1.74	0.88
1:A:191:PRO:HG3	1:A:247:THR:HG23	1.55	0.87
1:B:251:ASN:ND2	1:B:259:ARG:HE	1.73	0.87
1:B:177:SER:HA	1:B:282:ASN:ND2	1.93	0.83
1:A:58:SER:HB2	1:A:140:HIS:CE1	2.13	0.83
1:A:119:THR:OG1	1:A:122:GLU:HG3	1.78	0.82
1:A:142:ILE:HG23	1:A:171:ILE:HD12	1.59	0.82
1:B:268:GLU:CG	1:B:269:PRO:HD2	2.10	0.80
1:B:119:THR:OG1	1:B:122:GLU:HG3	1.82	0.79
1:A:177:SER:HA	1:A:282:ASN:ND2	1.96	0.78
1:B:58:SER:HB2	1:B:140:HIS:CE1	2.18	0.78
1:A:151:ILE:HB	1:A:152:PRO:HD3	1.63	0.78
1:B:251:ASN:HD22	1:B:259:ARG:HE	1.32	0.78
1:B:8:LEU:HA	1:B:248:CYS:SG	2.24	0.77
1:A:268:GLU:CG	1:A:269:PRO:HD2	2.16	0.76
1:B:151:ILE:HB	1:B:152:PRO:HD3	1.67	0.75
1:A:34:GLU:HG2	1:A:38:LYS:HD3	1.70	0.74
1:B:34:GLU:HG2	1:B:38:LYS:HD3	1.71	0.73
1:B:71:LEU:HD21	1:B:85:ASN:HD22	1.54	0.72
1:A:204:LEU:HD23	1:A:288:ILE:HD12	1.71	0.72
1:B:234:ASN:HD22	1:B:237:GLY:N	1.84	0.70
1:B:204:LEU:HD23	1:B:288:ILE:HD12	1.74	0.69
1:B:234:ASN:ND2	1:B:237:GLY:H	1.83	0.69
1:B:21:ARG:HG3	1:B:239:LEU:HD13	1.78	0.66
1:A:210:ASN:OD1	1:B:126:ARG:NH2	2.29	0.65
1:A:219:ASP:O	1:A:220:LEU:HD23	1.97	0.64
1:A:234:ASN:ND2	1:A:237:GLY:H	1.86	0.64
1:B:268:GLU:HG3	1:B:269:PRO:CD	2.25	0.64
1:B:234:ASN:HD21	1:B:236:ASN:HB3	1.62	0.64
1:B:219:ASP:O	1:B:220:LEU:HD23	1.97	0.64
1:B:275:LYS:O	1:B:277:GLY:N	2.31	0.64
1:A:203:ASP:O	1:A:207:MET:HG3	1.98	0.64
1:A:33:GLN:NE2	1:A:69:GLN:NE2	2.45	0.63
1:A:71:LEU:HD21	1:A:85:ASN:HD22	1.62	0.63

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:259:ARG:HG2	1:B:259:ARG:HH11	1.63	0.63
1:A:228:ASP:O	1:A:237:GLY:HA3	1.98	0.62
1:A:21:ARG:HG3	1:A:239:LEU:HD13	1.81	0.61
1:A:234:ASN:HD21	1:A:236:ASN:HB3	1.65	0.61
1:B:270:GLU:HG3	1:B:271:PHE:CD1	2.35	0.61
1:A:259:ARG:HG2	1:A:259:ARG:HH11	1.66	0.61
1:A:270:GLU:HG3	1:A:271:PHE:CD1	2.35	0.60
1:B:33:GLN:NE2	1:B:69:GLN:NE2	2.49	0.60
1:B:270:GLU:HG3	1:B:271:PHE:HD1	1.66	0.60
1:B:234:ASN:ND2	1:B:236:ASN:HB3	2.17	0.60
1:A:275:LYS:O	1:A:277:GLY:N	2.34	0.59
1:B:228:ASP:O	1:B:237:GLY:HA3	2.01	0.59
1:B:203:ASP:O	1:B:207:MET:HG3	2.01	0.59
1:B:259:ARG:HG2	1:B:259:ARG:NH1	2.17	0.59
1:B:93:GLU:O	1:B:97:LYS:HG3	2.03	0.59
1:B:166:ALA:O	1:B:167:LYS:HG2	2.03	0.59
1:A:234:ASN:HD22	1:A:237:GLY:N	1.86	0.58
1:A:168:MET:HE3	1:A:290:ILE:HD12	1.84	0.58
1:A:270:GLU:HG3	1:A:271:PHE:HD1	1.67	0.58
1:A:268:GLU:HG3	1:A:269:PRO:CD	2.29	0.58
1:A:36:MET:SD	1:A:40:LEU:HD22	2.44	0.58
1:B:19:PHE:HE2	1:B:23:LEU:HD11	1.69	0.58
1:A:234:ASN:ND2	1:A:236:ASN:HB3	2.19	0.57
1:A:86:GLU:HG3	1:A:136:TRP:HH2	1.69	0.57
1:A:259:ARG:HG2	1:A:259:ARG:NH1	2.19	0.56
1:B:21:ARG:CG	1:B:239:LEU:HD13	2.35	0.56
1:B:65:GLU:H	1:B:65:GLU:CD	2.09	0.56
1:B:261:GLU:O	1:B:265:ASP:HB2	2.06	0.56
1:A:19:PHE:HE2	1:A:23:LEU:HD11	1.71	0.56
1:B:19:PHE:CE2	1:B:23:LEU:HD11	2.41	0.56
1:A:166:ALA:O	1:A:167:LYS:HG2	2.06	0.55
1:A:193:ASP:OD1	1:A:195:LEU:HB2	2.06	0.55
1:B:197:GLN:HG2	1:B:199:ILE:HD13	1.90	0.54
1:A:21:ARG:CG	1:A:239:LEU:HD13	2.36	0.54
1:A:65:GLU:CD	1:A:65:GLU:H	2.11	0.54
1:A:197:GLN:HG2	1:A:199:ILE:HD13	1.90	0.54
1:B:32:MET:O	1:B:36:MET:HG2	2.08	0.53
1:A:78:TYR:HB3	1:A:81:VAL:HG21	1.90	0.53
1:B:193:ASP:OD1	1:B:195:LEU:HB2	2.09	0.53
1:B:71:LEU:HD21	1:B:85:ASN:ND2	2.24	0.53
1:A:151:ILE:HB	1:A:152:PRO:CD	2.38	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:19:PHE:CE2	1:A:23:LEU:HD11	2.44	0.52
1:A:261:GLU:O	1:A:265:ASP:HB2	2.09	0.52
1:B:193:ASP:OD1	1:B:194:ASP:N	2.42	0.52
1:B:21:ARG:NH1	1:B:236:ASN:OD1	2.43	0.52
1:B:86:GLU:HG3	1:B:136:TRP:HH2	1.74	0.52
1:A:33:GLN:HE22	1:A:69:GLN:CD	2.14	0.51
1:B:168:MET:HE3	1:B:290:ILE:HD12	1.92	0.50
1:A:193:ASP:OD1	1:A:194:ASP:N	2.44	0.50
1:A:32:MET:O	1:A:36:MET:HG2	2.12	0.50
1:B:17:GLU:O	1:B:21:ARG:HG2	2.11	0.50
1:A:168:MET:CE	1:A:290:ILE:HD12	2.41	0.50
1:A:193:ASP:OD1	1:A:195:LEU:N	2.24	0.50
1:B:19:PHE:O	1:B:22:PHE:HB3	2.12	0.50
1:B:276:GLU:C	1:B:278:LYS:H	2.14	0.50
1:B:35:PHE:HD2	1:B:36:MET:CE	2.25	0.49
1:A:89:GLU:OE2	1:A:90:PRO:HD2	2.12	0.49
1:B:262:LEU:O	1:B:263:GLY:C	2.51	0.48
1:A:33:GLN:CD	1:A:69:GLN:HE22	2.16	0.48
1:B:33:GLN:CD	1:B:69:GLN:HE22	2.17	0.48
1:B:36:MET:SD	1:B:40:LEU:HD22	2.54	0.48
1:B:228:ASP:HB3	1:B:234:ASN:ND2	2.28	0.48
1:A:33:GLN:OE1	1:A:69:GLN:OE1	2.31	0.48
1:B:63:ALA:HA	1:B:65:GLU:OE2	2.12	0.48
1:A:63:ALA:HA	1:A:65:GLU:OE2	2.13	0.48
1:A:33:GLN:HE22	1:A:69:GLN:NE2	2.12	0.47
1:A:93:GLU:O	1:A:97:LYS:HG3	2.14	0.47
1:A:269:PRO:HA	1:A:272:SER:O	2.14	0.47
1:B:33:GLN:OE1	1:B:69:GLN:OE1	2.32	0.47
1:A:21:ARG:NH1	1:A:236:ASN:OD1	2.47	0.47
1:B:139:ILE:HB	1:B:168:MET:HG3	1.97	0.47
1:B:269:PRO:HA	1:B:272:SER:O	2.15	0.47
1:A:222:SER:HB3	1:A:283:ASN:O	2.15	0.47
1:B:107:ASN:O	1:B:108:LEU:HD23	2.14	0.47
1:B:33:GLN:HE22	1:B:69:GLN:CD	2.17	0.47
1:B:168:MET:CE	1:B:290:ILE:HD12	2.43	0.47
1:A:46:ARG:HA	1:A:78:TYR:OH	2.14	0.47
1:B:89:GLU:OE2	1:B:90:PRO:HD2	2.15	0.47
1:A:17:GLU:O	1:A:21:ARG:HG2	2.14	0.47
1:A:276:GLU:C	1:A:278:LYS:H	2.18	0.47
1:B:9:PHE:HD2	1:B:195:LEU:HD23	1.80	0.47
1:A:228:ASP:HB3	1:A:234:ASN:ND2	2.30	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:8:LEU:HG	1:B:15:TYR:HD1	1.80	0.46
1:A:239:LEU:O	1:A:242:ASP:HB2	2.16	0.46
1:B:78:TYR:HB3	1:B:81:VAL:HG21	1.98	0.46
1:B:88:VAL:O	1:B:88:VAL:HG12	2.15	0.46
1:B:193:ASP:OD1	1:B:195:LEU:N	2.26	0.46
1:A:262:LEU:O	1:A:263:GLY:C	2.55	0.46
1:A:9:PHE:HD2	1:A:195:LEU:HD23	1.81	0.45
1:A:151:ILE:HG21	1:A:207:MET:CE	2.46	0.45
1:B:226:ILE:O	1:B:229:CYS:HB2	2.16	0.45
1:A:35:PHE:HD2	1:A:36:MET:CE	2.28	0.45
1:B:151:ILE:HG21	1:B:207:MET:CE	2.45	0.45
1:B:222:SER:HB3	1:B:283:ASN:O	2.16	0.45
1:A:19:PHE:O	1:A:22:PHE:HB3	2.17	0.45
1:A:58:SER:HB2	1:A:140:HIS:NE2	2.30	0.45
1:B:16:VAL:O	1:B:17:GLU:C	2.55	0.45
1:A:164:THR:HG22	1:A:165:ASN:OD1	2.17	0.44
1:A:28:GLU:O	1:A:31:CYS:HB3	2.17	0.44
1:B:34:GLU:CG	1:B:38:LYS:HD3	2.45	0.44
1:A:16:VAL:O	1:A:17:GLU:C	2.54	0.44
1:A:34:GLU:CG	1:A:38:LYS:HD3	2.43	0.44
1:A:89:GLU:O	1:A:117:LYS:HA	2.17	0.44
1:B:49:ASP:C	1:B:50:THR:HG23	2.37	0.44
1:A:249:ASN:O	1:A:250:PHE:C	2.56	0.44
1:B:33:GLN:HE22	1:B:69:GLN:NE2	2.15	0.44
1:B:231:ILE:O	1:B:232:ASP:C	2.56	0.44
1:B:100:GLU:O	1:B:104:LYS:HG3	2.18	0.44
1:B:46:ARG:HA	1:B:78:TYR:OH	2.18	0.43
1:B:276:GLU:C	1:B:278:LYS:N	2.71	0.43
1:B:44:ILE:HD12	1:B:44:ILE:O	2.18	0.43
1:A:8:LEU:HG	1:A:15:TYR:HD1	1.82	0.43
1:B:21:ARG:HG3	1:B:239:LEU:CD1	2.48	0.43
1:B:28:GLU:O	1:B:31:CYS:HB3	2.18	0.43
1:B:89:GLU:O	1:B:117:LYS:HA	2.18	0.43
1:B:151:ILE:HB	1:B:152:PRO:CD	2.42	0.43
1:A:231:ILE:O	1:A:232:ASP:C	2.58	0.42
1:A:18:SER:O	1:A:19:PHE:C	2.57	0.42
1:A:71:LEU:HD21	1:A:85:ASN:ND2	2.32	0.42
1:A:100:GLU:O	1:A:104:LYS:HG3	2.19	0.42
1:B:262:LEU:HD23	1:B:262:LEU:HA	1.82	0.42
1:A:8:LEU:HA	1:A:248:CYS:SG	2.60	0.42
1:A:18:SER:O	1:A:21:ARG:N	2.52	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:14:LYS:O	1:B:15:TYR:C	2.57	0.42
1:B:18:SER:O	1:B:19:PHE:C	2.58	0.42
1:B:239:LEU:O	1:B:242:ASP:HB2	2.20	0.42
1:A:5:MET:HE3	1:A:254:ALA:HB2	2.01	0.42
1:A:48:GLY:O	1:A:49:ASP:C	2.57	0.42
1:B:221:LEU:HD12	1:B:222:SER:H	1.84	0.42
1:A:30:GLN:O	1:A:33:GLN:HB3	2.20	0.42
1:B:94:GLN:O	1:B:95:ILE:C	2.56	0.42
1:B:258:LEU:C	1:B:260:ALA:N	2.73	0.42
1:B:177:SER:HA	1:B:282:ASN:HD21	1.82	0.41
1:A:33:GLN:CD	1:A:69:GLN:NE2	2.73	0.41
1:A:49:ASP:C	1:A:50:THR:HG23	2.41	0.41
1:A:98:TYR:O	1:A:101:LEU:HB3	2.21	0.41
1:B:274:LYS:HA	1:B:278:LYS:O	2.20	0.41
1:A:258:LEU:C	1:A:260:ALA:N	2.74	0.41
1:B:35:PHE:HD2	1:B:36:MET:HE2	1.85	0.41
1:A:40:LEU:HB3	1:A:41:PRO:HD3	2.03	0.41
1:A:231:ILE:O	1:A:231:ILE:HG22	2.21	0.41
1:B:169:LEU:HD12	1:B:288:ILE:O	2.21	0.41
1:A:44:ILE:HD12	1:A:44:ILE:O	2.21	0.41
1:B:18:SER:O	1:B:21:ARG:N	2.54	0.41
1:A:139:ILE:HB	1:A:168:MET:HG3	2.03	0.40
1:B:128:LEU:HD13	1:B:128:LEU:HA	1.87	0.40
1:B:196:CYS:SG	1:B:197:GLN:N	2.94	0.40
1:B:241:TRP:HZ3	1:B:262:LEU:HD13	1.86	0.40
1:A:269:PRO:O	1:A:271:PHE:N	2.54	0.40
1:A:276:GLU:C	1:A:278:LYS:N	2.75	0.40
1:A:94:GLN:O	1:A:95:ILE:C	2.60	0.40
1:A:110:ASN:N	1:A:110:ASN:HD22	2.19	0.40
1:B:9:PHE:CD2	1:B:195:LEU:HD23	2.56	0.40

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:24:ASN:OD1	1:A:24:ASN:OD1[4_655]	2.19	0.01

5.3 Torsion angles ⓘ

5.3.1 Protein backbone ⓘ

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	286/292 (98%)	242 (85%)	37 (13%)	7 (2%)	6	26
1	B	286/292 (98%)	243 (85%)	39 (14%)	4 (1%)	11	39
All	All	572/584 (98%)	485 (85%)	76 (13%)	11 (2%)	8	32

All (11) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	249	ASN
1	A	276	GLU
1	B	276	GLU
1	A	250	PHE
1	A	270	GLU
1	B	12	HIS
1	A	12	HIS
1	A	15	TYR
1	B	15	TYR
1	A	187	GLY
1	B	187	GLY

5.3.2 Protein sidechains ⓘ

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the sidechain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	238/260 (92%)	225 (94%)	13 (6%)	21	52
1	B	238/260 (92%)	224 (94%)	14 (6%)	19	50

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
All	All	476/520 (92%)	449 (94%)	27 (6%)	20	51

All (27) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	30	GLN
1	A	65	GLU
1	A	67	ASP
1	A	82	CYS
1	A	107	ASN
1	A	154	THR
1	A	194	ASP
1	A	195	LEU
1	A	199	ILE
1	A	202	ASP
1	A	246	GLU
1	A	247	THR
1	A	265	ASP
1	B	30	GLN
1	B	65	GLU
1	B	67	ASP
1	B	82	CYS
1	B	107	ASN
1	B	154	THR
1	B	194	ASP
1	B	195	LEU
1	B	199	ILE
1	B	202	ASP
1	B	204	LEU
1	B	246	GLU
1	B	247	THR
1	B	265	ASP

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (20) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	29	HIS
1	A	30	GLN
1	A	69	GLN
1	A	84	ASN
1	A	85	ASN

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Mol	Chain	Res	Type
1	A	107	ASN
1	A	110	ASN
1	A	197	GLN
1	A	206	GLN
1	A	234	ASN
1	B	30	GLN
1	B	69	GLN
1	B	84	ASN
1	B	85	ASN
1	B	110	ASN
1	B	124	GLN
1	B	197	GLN
1	B	206	GLN
1	B	234	ASN
1	B	251	ASN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

There are no non-standard protein/DNA/RNA residues in this entry.

5.5 Carbohydrates ⓘ

There are no monosaccharides in this entry.

5.6 Ligand geometry ⓘ

2 ligands are modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	#|Z| > 2	Counts	RMSZ	#|Z| > 2
2	THA	A	400	-	17,17,17	6.15	10 (58%)	21,24,24	2.95	11 (52%)
2	THA	B	401	-	17,17,17	5.59	11 (64%)	21,24,24	2.73	13 (61%)

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	THA	A	400	-	-	-	0/3/3/3
2	THA	B	401	-	-	-	0/3/3/3

All (21) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	400	THA	C8-N7	14.99	1.51	1.32
2	B	401	THA	C8-N7	13.10	1.49	1.32
2	A	400	THA	C1-C2	11.30	1.62	1.36
2	B	401	THA	C1-C2	10.03	1.59	1.36
2	A	400	THA	C6-C5	8.47	1.56	1.36
2	B	401	THA	C6-C5	7.68	1.54	1.36
2	A	400	THA	C9-C8	7.06	1.51	1.40
2	A	400	THA	C6-C1	6.74	1.55	1.38
2	B	401	THA	C9-C8	6.62	1.50	1.40
2	B	401	THA	C5-C4	5.70	1.53	1.42
2	B	401	THA	C6-C1	5.68	1.53	1.38
2	A	400	THA	C5-C4	5.50	1.53	1.42
2	A	400	THA	C2-C3	5.19	1.50	1.41
2	B	401	THA	C14-C9	-5.00	1.43	1.51
2	A	400	THA	C14-C9	-4.74	1.44	1.51
2	B	401	THA	C2-C3	4.52	1.49	1.41
2	B	401	THA	C10-N15	4.28	1.48	1.36
2	A	400	THA	C10-N15	4.17	1.48	1.36
2	B	401	THA	C4-C3	3.83	1.48	1.42
2	A	400	THA	C4-C3	3.50	1.48	1.42
2	B	401	THA	C11-C8	-2.51	1.46	1.50

All (24) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	400	THA	C9-C8-N7	-7.30	117.44	123.68
2	B	401	THA	C9-C8-N7	-6.55	118.08	123.68

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
2	B	401	THA	C1-C6-C5	-4.17	114.60	120.44
2	B	401	THA	C1-C2-C3	-3.82	114.58	120.08
2	A	400	THA	C1-C2-C3	-3.78	114.64	120.08
2	A	400	THA	C11-C8-N7	3.77	122.79	116.73
2	A	400	THA	C1-C6-C5	-3.74	115.20	120.44
2	A	400	THA	C12-C11-C8	3.71	119.63	113.53
2	A	400	THA	C5-C4-C3	3.66	122.42	118.33
2	B	401	THA	C8-N7-C3	3.44	121.91	117.67
2	B	401	THA	C12-C11-C8	3.38	119.08	113.53
2	A	400	THA	C8-N7-C3	3.34	121.79	117.67
2	B	401	THA	C13-C14-C9	3.31	119.63	112.84
2	B	401	THA	C5-C4-C3	3.16	121.86	118.33
2	A	400	THA	C13-C14-C9	2.98	118.97	112.84
2	B	401	THA	C11-C8-N7	2.97	121.51	116.73
2	A	400	THA	C2-C3-C4	2.79	122.33	119.13
2	A	400	THA	C5-C4-C10	-2.61	117.40	123.08
2	A	400	THA	C12-C13-C14	2.40	123.19	112.55
2	B	401	THA	C5-C4-C10	-2.27	118.15	123.08
2	B	401	THA	C4-C3-N7	-2.25	120.43	122.81
2	B	401	THA	C2-C3-C4	2.16	121.61	119.13
2	B	401	THA	C12-C13-C14	2.13	122.00	112.55
2	B	401	THA	C6-C5-C4	-2.02	118.10	120.89

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

5.7 Other polymers ⓘ

There are no such residues in this entry.

5.8 Polymer linkage issues ⓘ

There are no chain breaks in this entry.

Full wwPDB X-ray Structure Validation Report

6 Fit of model and data ⓘ

6.1 Protein, DNA and RNA chains ⓘ

EDS was not executed - this section is therefore empty.

6.2 Non-standard residues in protein, DNA, RNA chains ⓘ

EDS was not executed - this section is therefore empty.

6.3 Carbohydrates ⓘ

EDS was not executed - this section is therefore empty.

6.4 Ligands ⓘ

EDS was not executed - this section is therefore empty.

6.5 Other polymers ⓘ

EDS was not executed - this section is therefore empty.